AMENDMENTS TO THE CLAIMS:

This listing of claims will replace all prior versions, and listings, of claims in the application:

1 (currently amended). A compound of formula I,

$$R^1$$
 P^2
 P^2
 P^3
 P^2
 P^3
 P^3

wherein

one of R1 and R2 represents a structural fragment of formula la

and the other represents R4;

Z represents O or N(R⁵);

 R^3 represents one or more optional substituents selected from OH, halo, cyano, nitro, $C(O)OR^6$, C_{1-6} alkoxy or C_{1-6} alkyl, (which two latter groups are optionally substituted and/or terminated by one or more halo or hydroxy group), or $N(R^7)R^8$;

R⁴ represents H, OH, halo, cyano, nitro, C(O)OR⁶, C₁₋₆ alkoxy or C₁₋₆ alkyl, (which two latter groups are optionally substituted and/or terminated by one or more halo or hydroxy group), or N(R⁷)R⁸;

Ar¹ represents phenyl, C_{1-3} alkylphenyl, C_{1-3} alkyldiphenyl, C_{3-7} cycloalkyl, C_{1-3} -alkyl- C_{3-7} -cycloalkyl, C_{1-3} -alkyl-di- C_{3-7} -cycloalkyl, naphthyl, C_{1-3} alkylnaphthyl, thienyl, imidazolyl or isoxazolyl, all of which may be substituted by one or more substituent selected from OH, halo, cyano, nitro, $C(O)OR^6$, C_{1-6} alkoxy or C_{1-6} alkyl, (which two latter groups are optionally substituted and/or terminated by one or more halo or hydroxy group), or $N(R^7)R^8$;

 R^5 represents H, C_{1-6} alkyl, phenyl or C_{1-3} alkylphenyl, (which three latter groups are optionally substituted and/or terminated by one or more substituent selected from OH, halo, cyano, nitro, $C(O)OR^9$, $C(O)N(R^{10})R^{11}$, $P(O)(R^{12})R^{13}$, $P(O)(OR^{14})OR^{15}$, $S(O)_2(R^{16})R^{17}$, $S(O)_2N(R^{18})R^{19}$, C_{1-6} alkoxy or C_{1-6} alkyl, (which two latter groups are optionally substituted and/or terminated by one or more halo or hydroxy group) or $N(R^{20})R^{21}$;

Y represents O, S, S(O), S(O)₂ or $N(R^{22})$;

 R^{10} and R^{11} independently represent H, OR^{23} , $C(O)R^{24}$, $OC(O)R^{25}$, $C(O)OR^{26}$, C_{1-4} alkyl, (which latter group is optionally substituted and/or terminated by one or more substituent selected from C_{1-4} alkyl, OR^{27} , $N(R^{28})R^{29}$, $C(O)OR^{30}$, $C(O)N(R^{31})R^{32}$, $P(O)(R^{33})R^{34}$, $P(O)(OR^{35})OR^{36}$ and $S(O)_2N(R^{37})R^{38}$), $-(CH_2CH_2O-)_pR^{39}$ or, together with the nitrogen atom to which they are attached, form a C_{4-7} nitrogen-containing, aromatic or non-aromatic, ring which ring may contain a further heteroatom or group (as appropriate) selected from O, S and $N(R^{40})$ and may further be substituted by one or

more substituent selected from C(O)R⁴¹, C(O)OR⁴² or C(O)N(R⁴³)R⁴⁴;

 R^{28} , R^{29} , R^{30} , R^{31} , R^{32} and R^{40} independently represent H or C_{1-6} alkyl, which latter group is optionally substituted and/or terminated by one or more substituent selected from $C(O)R^{45}$, $C(O)OR^{46}$ or $C(O)N(R^{47})R^{48}$;

at each occurrence, R⁶, R⁷ and R⁸ independently represent H or C₁₋₄ alkyl;

 R^9 , R^{12} , R^{13} , R^{14} , R^{15} , R^{16} , R^{17} , R^{18} , R^{19} , R^{20} , R^{21} , R^{22} , R^{23} , R^{24} , R^{25} , R^{26} , R^{27} , R^{33} , R^{34} , R^{35} , R^{36} , R^{37} , R^{38} , R^{39} , R^{41} , R^{42} , R^{43} , R^{44} , R^{45} , R^{46} , R^{47} and R^{48} independently represent H or C_{1-4} alkyl;

n represents 2;

p represents 1, 2, 3, 4, 5 or 6; and

B represents a structural fragment of formula lb, lc, ld or le

wherein

X¹ and X² independently represent a single bond or CH₂; or a pharmaceutically acceptable salt thereof.

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2 (original). A compound of formula I, as defined in claim 1, wherein, when B represents a structural fragment of formula Ib, Id, Ie or Ic in which latter fragment X^1 and X^2 both represent CH_2 , then n represents 2.

3 (cancelled).

4 (previously presented). A compound of formula I, as defined in claim 1, wherein R² represents a structural fragment of formula Ia and R¹ represents R⁴.

5 (previously presented). A compound of formula I, as defined in claim 1, wherein Z represents O or $N(R^5)$, in which latter case R^5 represents C_{1-6} alkyl terminated by $C(O)N(R^{10})R^{11}$.

6 (previously presented). A compound of formula I, as defined in claim 1, wherein R³ is not present, or represents methyl, chloro or methoxy.

7 (previously presented). A compound of formula I, as defined in claim 1, wherein Ar¹ represents substituted phenyl.

8 (previously presented). A compound of formula I, as defined in claim 1 wherein Y represents O.

9 (previously presented). A compound of formula I, as defined in claim 1

wherein B represents a structural fragment of formula lb.

10 (original). A compound as claimed in Claim 1 which is:

N-{3-[2-(4-aminoiminomethylphenyl)ethoxy]phenyl} benzenesulfonamide;

benzenesulfonic acid- {3-[2-(4-aminoiminomethylphenyl)ethoxy]-5-methyl}phenyl ester;

N-{3-[2-(4-aminoiminomethylphenyl)ethoxy]phenyl}-2-chlorobenzenesulfonamide;

N-{3-[2-(4-aminoiminomethylphenyl)ethoxy]phenyl}-2-cyanobenzene-

sulfonamide;

N-{3-[2-(4-aminoiminomethylphenyl)ethoxy]phenyl}-2-fluorobenzene-sulfonamide:

N-{3-[2-(4-aminoiminomethylphenyl)ethoxy]phenyl}-2-(trifluoromethoxy)-benzenesulfonamide;

N-{3-[2-(4- aminoiminomethylphenyl)ethoxy]phenyl}-4-fluorobenzene-sulfonamide;

N-{3-[2-(4-aminoiminomethylphenyl)ethoxy]phenyl}-2,5-dimethylbenzene-sulfonamide:

N-{3-[2-(4-aminoiminomethylphenyl)ethoxy]phenyl}-5-chlorothiophene-2-sulfonamide;

N-{3-[2-(4-aminoiminomethylphenyl)ethoxy]phenyl}-1-methylimidazole-3-sulfonamide;

N-{3-[2-(4-aminoiminomethylphenyl)ethoxy]phenyl}-3,5-dimethylisoxazole-4-sulfonamide;

N-{3-[2-(4-aminoiminomethylphenyl)ethoxy]phenyl} benzylsulfonamide;

N-{3-[2-(4-aminoiminomethylphenyl)ethoxy]phenyl}-2,5-dichiorothiophene-3-sulfonamide;

N-{3-[2-(4-aminoiminomethylphenyl)ethoxy]-5-methylphenyl}-2-chlorobenzenesulfonamide;

N-{3-[2-(4-aminoiminomethylphenyl)ethoxy]-2-methylphenyl}-benzenesulfonamide;

N-{5-[2-(4-aminoiminomethylphenyl)ethoxy]-2-methylphenyl}benzenesulfonamide;

N-{3-[2-(4-aminoiminomethylphenyl)ethoxy]-5-methylphenyl} benzenesulfonamide;

N-{3-[2-(4-aminoiminomethylphenyl)ethylthio]phenyl} benzenesulfonamide;

N-(2-chlorophenyl)sulfonyl-3-[2-(4-aminoiminomethylphenyl) ethoxy]-5-methylphenylaminoacetic acid, ethyl ester;

N-(2-chlorophenyl)sulfonyl-3-[2-(4-aminoiminomethylphenyl)ethoxy]-5-methylphenylaminoacetamide;

N-(2-chlorophenyl)sulfonyl-3-[2-(4-aminoiminomethylphenyl)ethoxy]-5-methylphenylaminoacetic acid;

N-(2-chlorophenyl)sulfonyl-2-{3-[2-(4-aminoiminomethylphenyl)ethoxy]-5-methylphenylamino}propanoic acid, ethyl ester;

2-{3-[2-(4-aminoiminomethylphenyl)ethoxy]-N-(2-chlorophenyl)sulfonyl-5-methylphenylamino}propanamide;

N-(2-chlorophenyl)sulfonyl-2-{3-[2-(4-aminoiminomethylphenyl)ethoxy]-5-methylphenylamino}propanoic acid;

N-(2-chlorophenyl)sulfonyl-2-{3-[2-(4-aminoiminomethylphenyl)ethoxy]-5-methylphenylamino}propanoic acid, methyl ester;

N-(2-chlorophenyl)sulfonyl-3-{3-[2-(4-aminoiminomethylphenyl)ethoxy]-5-methylphenylamino}butanoic acid, ethyl ester;

3-{3-[2-(4-aminoiminomethylphenyl)ethoxy]-N-(2-chlorophenyl)sulfonyl-5-methylphenylamino}butanamide;

N-(2-chlorophenyl)sulfonyl-3-{3-[2-(4-aminoiminomethylphenyl)ethoxy]-5-methylphenylamino}butanoic acid;

N-(2-chlorophenyl)sulfonyl-4-{3-[2-(4-aminoiminomethylphenyl)ethoxy]-5-methylphenylamino}pentanoic acid, ethyl ester;

4-{3-[2-(4-aminoiminomethylpheny)ethoxy]-N-(2-chlorophenyl)sulfonyl-5-methylphenylamino}pentanamide;

N-(2-chlorophenyl)sulfonyl-4-{3-[2-(4-aminoiminomethylphenyl)ethoxy]-5-methylphenylamino}pentanoic acid;

N-(2-chlorophenyl)sulfonyl-5-{3-[2-(4-aminoiminomethylphenyl)ethoxy]-5-methylphenylamino}hexanoic acid, ethyl ester;

5-{3-[2-(4-aminoiminomethylphenyl)ethoxy]- N-(2-chlorophenyl)sulfonyl-5-methylphenyl amino}pentanamide;

N-(2-chlorophenyl)sulfonyl-5-{3-[2-(4-aminoiminomethylphenyl)ethoxy]-5-methylphenylamino}hexanoic acid;

N-phenylsulfonyl-3-[2-(4-aminoiminomethylphenyl)ethoxy]phenylaminoacetic acid, ethyl ester;

N-phenylsulfonyl-3-[2-(4- aminoiminomethylphenyl)ethoxy]phenylaminoacetic

acid;

N-{3-[2-(4-aminoiminomethylphenyl)ethoxy]phenyl}-N-(2-hydroxyethyl)-benzenesulfonamide;

N-{3-[2-(4-aminoiminomethylphenyl)ethoxy]phenyl}-N-(dimethyloxophosphinylmethyl)-benzenesulfonamide;

2-chlorobenzenesulfonic acid, 3-[2-(4-aminoiminomethylphenyl)ethoxy]-5-methylphenyl ester;

benzenesulfonic acid, 3-[2-(4-aminoiminomethylphenyl)ethoxy]phenyl ester;

2-chloro-4-fluorobenzenesulfonic acid, 3-[2-(4-aminoiminomethylphenyl)-ethoxy]-5-chlorophenyl ester;

2-chlorobenzenesulfonic acid, 3-[2-(4-aminoiminomethylphenyl)ethoxy]-5-methoxyphenyl ester;

2-chlorobenzenesulfonic acid, 3-[2-(4-aminoiminomethylphenyl)ethoxy]-5ethylphenyl ester;

N-{2-[2-(4-aminoiminomethylphenyl)ethylthio]phenyl}benzenesulfonamide;

N-{2-[2-(4-aminoiminomethylphenyl)ethylthio]phenyl}-2,4,5-trichlorobenzenesulfonamide;

N-{2-[2-(4-aminoiminomethylphenyl)ethylthio]phenyl}-2-chloro-5-methoxybenzenesulfonamide;

N-{2-[2-(4-aminoiminomethylphenyl)ethylthio]phenyl}-2,5-dibromobenzenesulfonamide;

N-{2-[2-(4-aminoiminomethylphenyl)ethylthio]phenyl}-2,5-dichlorobenzenesulfonamide; N-{2-[2-(4-aminoiminomethylphenyl)-ethylthio]-phenyl}-2-methoxy-5-methylbenzenesulfonamide;

N-{2-[2-(4-aminoiminomethylphenyl)ethylthio]phenyl}-2,3,5,6-tetramethylbemzenesulfonamide;

N-{2-[2-(4-aminoiminomethylphenyl)ethylthio]phenyl}-3,4-dimethoxy-benzenesulfonamide;

N-{2-[2-(4-aminoiminomethylphenyl)ethylthio]phenyl}-3-bromobenzenesulfonamide;

N-{2-[2-(4-aminoiminomethylphenyl)ethylthio]phenyl}-3,4-dibromobenzene-sulfonamide;

N-{2-[2-(4-aminoiminomethylphenyl)ethylthio]phenyl}-2-chloro-4-fluorobenzenesulfonamide; or

N-{2-[2-(4-aminoiminomethylphenyl)ethylthio]phenyl}- 5-bromo-2-methoxybenzenesulfonamide.

11 (original). A compound of formula I, as defined in claim 1, provided that R¹ represents a structural fragment of formula Ia and R² represents R⁴.

12 (original). A compound of formula I, as defined in claim 1, provided that Ar¹ represents optionally substituted phenyl.

13 (original). A compound of formula I, as defined in claim 1, provided that R⁵ is not substituted by P(O)(OR¹⁴)OR¹⁵, S(O)₂(R¹⁶)R¹⁷ or S(O)₂N(R¹⁸)R¹⁹.

14 (original). A compound of formula I, as defined in claim 1, provided that R^{10} and/or R^{11} represent H or unsubstituted C_{1-4} alkyl.

15 (original). A compound of formula I, as defined in claim 1, provided that Y represents O, S or N(R⁵).

16 (original). A compound of formula I, as defined in claim 1, provided that B represents a structural fragment of formula Ib, Ic, or Id.

17 (original). A compound of formula I, as defined in claim 1, provided that R² represents a structural fragment of formula Ia and R¹ represents R⁴.

18 (original). A compound of formula I, as defined in claim 1, provided that Ar¹ does not represent optionally substituted phenyl.

19 (original). A compound of formula I, as defined in claim 1, provided that R^5 is substituted by $P(O)(OR^{14})OR^{15}$, $S(O)_2(R^{16})R^{17}$ or $S(O)_2N(R^{18})R^{19}$.

20 (original). A compound of formula I, as defined in claim 1, provided that R¹⁰ and/or R¹¹ do not represent H or unsubstituted C₁₋₄ alkyl.

21 (original). A compound of formula I, as defined in claim 1, provided that Y

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represents S(O) or S(O)₂.

22 (original). A compound of formula I as defined in claim 1, provided that B represents a structural fragment of formula le.

23 (previously presented). A pharmaceutical formulation including a compound as defined in claim 1, or a pharmaceutically acceptable salt thereof, in admixture with a pharmaceutically acceptable adjuvant, diluent or carrier.

24-34 (canceled).